

STIC Search Report Biotech-Chem Library

STIC Database Tracking Number: 138289

TO: Andrew D Kosar

Location: rem/3c04/3c18

Art Unit: 1654

Wednesday, December 01, 2004

Case Serial Number: 10/613754

From: Peggy Ruppel

Location: Biotech-Chem Library

REMSEN 1B65

Phone: 571-272-2557

Peggy.Ruppel@uspto.gov

Search Notes

It was simpler for me to bundle the results of the two searches that you submitted for this application together, since the results sets were so small.

Please contact me if you have any questions or comments.

Thank you for using STIC services.



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STRUCTURE FILE UPDATES: 29 NOV 2004 HIGHEST RN 790629-40-2 DICTIONARY FILE UPDATES: 29 NOV 2004 HIGHEST RN 790629-40-2

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

NODE ATTRIBUTES: DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES: RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 13

STEREO ATTRIBUTES: NONE

L69 2556 SEA FILE=REGISTRY SSS FUL L65

=> d ide 169 1-7

L69 ANSWER 1 OF 2556 REGISTRY COPYRIGHT 2004 ACS on STN

RN 790602-32-3 REGISTRY

CN INDEX NAME NOT YET ASSIGNED

MF C19 H19 N5

CI COM

SR CA

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L69 ANSWER 2 OF 2556 REGISTRY COPYRIGHT 2004 ACS on STN

RN 790151-43-8 REGISTRY

CN 1H-Pyrazolo[3,4-b]quinoline-1-acetic acid, 4-[[3-(dimethylamino)propyl]amino]-3-methyl-, methyl ester (9CI) (CA INDEX NAME)

MF C19 H25 N5 O2

CI COM

SR CA

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L69 ANSWER 3 OF 2556 REGISTRY COPYRIGHT 2004 ACS on STN

RN 789481-53-4 REGISTRY

CN lH-Pyrazolo[3,4-b]quinolin-4-amine, 8-methoxy-3-methyl-1-(2-pyridinyl)-(9CI) (CA INDEX NAME)

MF C17 H15 N5 O

CI COM

SR CA

L69 ANSWER 4 OF 2556 REGISTRY COPYRIGHT 2004 ACS on STN

RN 789433-46-1 REGISTRY

CN 1,3-Propanediamine, N,N-dimethyl-N'-[3-methyl-1-(phenylmethyl)-1H-pyrazolo[3,4-b]quinolin-4-yl]- (9CI) (CA INDEX NAME)

MF C23 H27 N5

CI COM

SR CA

L69 ANSWER 5 OF 2556 REGISTRY COPYRIGHT 2004 ACS on STN

RN 788765-58-2 REGISTRY

CN 1-Propanamine, 3-[(1,3-dimethyl-1H-pyrazolo[3,4-b]quinolin-4-yl)thio]-N,N-dimethyl- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C17 H22 N4 S

CI COM

SR CA

L69 ANSWER 6 OF 2556 REGISTRY COPYRIGHT 2004 ACS on STN RN 788100-81-2 REGISTRY

CN 1,3-Propanediamine, N'-[1-[3-(dimethylamino)propyl]-3-methyl-1H-pyrazolo[3,4-b]quinolin-4-yl]-N,N-dimethyl- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C21 H32 N6

CI COM

SR CA

L69 ANSWER 7 OF 2556 REGISTRY COPYRIGHT 2004 ACS on STN

RN 787492-13-1 REGISTRY

CN 1,3-Propanediamine, N,N-dimethyl-N'-(1,3,5-trimethyl-1H-pyrazolo[3,4-b]quinolin-4-yl)- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C18 H25 N5

CI COM

SR CA

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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FILE 'HOME' ENTERED AT 15:34:36 ON 01 DEC 2004

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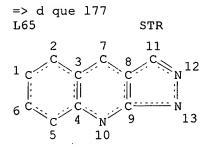
=> b hcaplus FILE 'HCAPLUS' ENTERED AT 15:51:52 ON 01 DEC 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 1 Dec 2004 VOL 141 ISS 23 FILE LAST UPDATED: 29 Nov 2004 (20041129/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

'OBI' IS DEFAULT SEARCH FIELD FOR 'HCAPLUS' FILE



NODE ATTRIBUTES: DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES: RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 13

STEREO ATTRIBUTES: NONE

L69 2556 SEA FILE=REGISTRY SSS FUL L65

L75 1292 SEA FILE=HCAPLUS ABB=ON PLU=ON "ERYTHROPOIETIN RECEPTORS"+OLD

/CT OR (EPO(A) RECEPT? OR EPOETIN(2A) RECEPT?)/BI

L76 52 SEA FILE=HCAPLUS ABB=ON PLU=ON L69(L)(USES+NT)/RL

L77 1 SEA FILE=HCAPLUS ABB=ON PLU=ON L76 AND L75

=> d ibib abs hitstr 177

L77 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2004 ACS on STN

ACCESSION NUMBER: 2004:41501 HCAPLUS

DOCUMENT NUMBER: 140:87744

TITLE: Affinity small molecules for the EPO

Searched by P. Ruppel

```
receptor
```

INVENTOR(S):

PATENT ASSIGNEE(S):

SOURCE:

Olsson, Lennart; Naranda, Tatjana

Receptron, Inc., USA PCT Int. Appl., 85 pp.

CODEN: PIXXD2

DOCUMENT TYPE:

LANGUAGE:

Patent English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO. KIND DATE APPLICATION NO. DATE ---------A2 WO 2004005323 20040115 WO 2003-US21394 20030703 WO 2004005323 A3 20040701 W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH,

PL, PT, RO, RU, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY,

KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

US 2004171541 A1 20040902 US 2003-613754 20030702 US 2004116346 A1 20040617 US 2003-612885 20030703 PRIORITY APPLN. INFO.: US 2002-393360P Ρ 20020703

US 2002-393361P Ρ 20020703 US 2002-394110P P 20020703

OTHER SOURCE(S): MARPAT 140:87744

Compds. are provided that complex with the modulating domain of erythropoietin receptor (EPO-R) for use with EPO-R to determine the presence of EPO-R, the ability of other mols. to bind to the modulating domain in competitive assays and to induce a signal by EPO-R into a cell when bound by the subject compds. in a physiol. environment. The compds. are characterized by having a six-membered heterocyclic ring comprising at least one nitrogen atom and include substituted triazolopyrimidine, pyridazinone, pyridine and piperidine.

IT645337-25-3

> RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)

(affinity small mols. for erythropoietin (EPO)

receptor and EPO receptor modulating

sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

645337-25-3 HCAPLUS RN

CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)

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=>

=> b reg FILE 'REGISTRY' ENTERED AT 15:35:00 ON 01 DEC 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 American Chemical Society (ACS)

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STRUCTURE FILE UPDATES: 29 NOV 2004 HIGHEST RN 790629-40-2 DICTIONARY FILE UPDATES: 29 NOV 2004 HIGHEST RN 790629-40-2

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

NODE ATTRIBUTES: DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES: RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 13

STEREO ATTRIBUTES: NONE L67 STR

NODE ATTRIBUTES:
DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS 22

STEREO ATTRIBUTES: NONE

L69 2556 SEA FILE=REGISTRY SSS FUL L65

L71 8 SEA FILE=REGISTRY SUB=L69 SSS FUL L67

=> d ide 171 1-8

L71 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

RN 748146-78-3 REGISTRY

CN INDEX NAME NOT YET ASSIGNED

FS 3D CONCORD

MF C33 H34 N4 O5 S

SR Chemical Library

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L71 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

RN 748146-41-0 REGISTRY

CN INDEX NAME NOT YET ASSIGNED

FS 3D CONCORD

MF C27 H32 N4 O4 S

SR Chemical Library

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L71 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

RN 748145-15-5 REGISTRY

CN INDEX NAME NOT YET ASSIGNED

FS 3D CONCORD

MF C25 H27 N3 O2

SR Chemical Library

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L71 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

RN 645337-25-3 REGISTRY

CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C17 H19 N3 O2

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

DT.CA CAplus document type: Patent

RL.P Roles from patents: BIOL (Biological study); USES (Uses)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)

1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L71 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

RN 521318-71-8 REGISTRY

CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 1-(4,6-dimethyl-2-pyrimidinyl)-

1,4,6,7,8,9-hexahydro-3,7,7-trimethyl-4-(5-methyl-2-furanyl)- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C24 H27 N5 O2

SR Chemical Library

LC STN Files: CHEMCATS

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L71 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

RN 521284-01-5 REGISTRY

CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 1-(4,6-dimethyl-2-pyrimidinyl)-4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C23 H25 N5 O2

SR Chemical Library

LC STN Files: CHEMCATS

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L71 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

RN. 380450-98-6 REGISTRY

CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl-1-(4-methylphenyl)- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C24 H25 N3 O2

SR Chemical Library

LC STN Files: CHEMCATS

L71 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

RN 378189-53-8 REGISTRY

CN Benzenesulfonamide, N,N-diethyl-4-[5-(4,5,6,7,8,9-hexahydro-3,7,7-trimethyl-5-oxo-1-phenyl-1H-pyrazolo[3,4-b]quinolin-4-yl)-2-furanyl]-(9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C33 H36 N4 O4 S

SR Chemical Library

LC STN Files: CHEMCATS

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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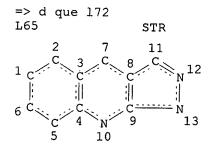
=> b hcaplus FILE 'HCAPLUS' ENTERED AT 15:39:03 ON 01 DEC 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 1 Dec 2004 VOL 141 ISS 23 FILE LAST UPDATED: 29 Nov 2004 (20041129/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

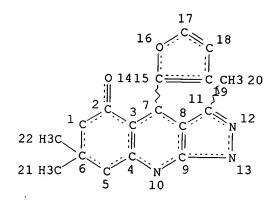
'OBI' IS DEFAULT SEARCH FIELD FOR 'HCAPLUS' FILE



NODE ATTRIBUTES: DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES: RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 13

STEREO ATTRIBUTES: NONE L67 STR



NODE ATTRIBUTES: DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES: RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 22

STEREO ATTRIBUTES: NONE

L69 2556 SEA FILE=REGISTRY SSS FUL L65

L71 8 SEA FILE=REGISTRY SUB=L69 SSS FUL L67 L72 1 SEA FILE=HCAPLUS ABB=ON PLU=ON L71

=> d ibib abs hitstr 172

L72 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2004 ACS on STN

ACCESSION NUMBER: 2004:41501 HCAPLUS

DOCUMENT NUMBER: 140:87744

TITLE:

Affinity small molecules for the EPO receptor

INVENTOR(S): Olsson, Lennart; Naranda, Tatjana

PATENT ASSIGNEE(S): Receptron, Inc., USA

SOURCE: PCT Int. Appl., 85 pp.

CODEN: PIXXD2 DOCUMENT TYPE:

Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.				KIND		DATE			APPLICATION NO.					DATE		
WO 2004005323 WO 2004005323					20040115			WO 2003-US21394					20030703			
	CO, GM, LS, PL, UG,	CR, HR, LT, PT, UZ,	CU, HU, LU, RO, VN,	CZ, ID, LV, RU, YU,	DE, IL, MA, SD, ZA,	AU, DK, IN, MD, SE, ZM,	DM, IS, MG, SG, ZW	DZ, JP, MK, SK,	EC, KE, MN, SL,	EE, KG, MW, TJ,	ES, KP, MX, TM,	FI, KR, MZ, TN,	GB, KZ, NO, TR,	GD, LC, NZ, TT,	GE, LK, OM, TZ,	GH, LR, PH, UA,
RW:	KG,	KZ,	MD,	RU,	ТJ,	MZ, TM, IE,	ΑT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK.	EE.	ES.

BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG US 2004171541 A1 20040902 US 2003-613754 20030702 US 2004116346 A 1 20040617 US 2003-612885 20030703 PRIORITY APPLN. INFO.: US 2002-393360P Ρ 20020703 US 2002-393361P Ρ 20020703 US 2002-394110P 20020703

OTHER SOURCE(S): MARPAT 140:87744

Compds. are provided that complex with the modulating domain of erythropoietin receptor (EPO-R) for use with EPO-R to determine the presence of EPO-R, the ability of other mols. to bind to the modulating domain in competitive assays and to induce a signal by EPO-R into a cell when bound by the subject compds. in a physiol. environment. The compds. are characterized by having a six-membered heterocyclic ring comprising at least one nitrogen atom and include substituted triazolopyrimidine, pyridazinone, pyridine and piperidine. ΙT

RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)

(affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

RN 645337-25-3 HCAPLUS

5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-CN 3,7,7-trimethyl- (9CI) (CA INDEX NAME)

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